



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 8, 2014 – 07:19 PM EDT

PDB ID : 4LP6
Title : Crystal Structure of Human Carbonic Anhydrase II in complex with a quino-
line oligoamide foldamer
Authors : Buratto, J.; Granier, T.; Langlois D'estaintot, B.; Huc, I.; Gallois, B.
Deposited on : 2013-07-15
Resolution : 2.15 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

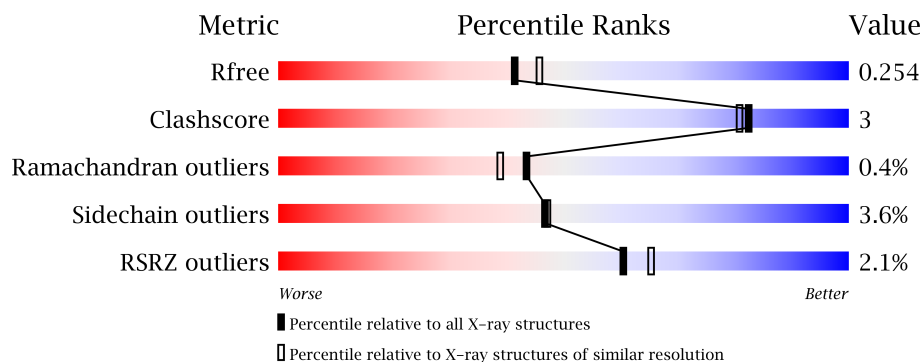
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23489
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23489

1 Overall quality at a glance



The reported resolution of this entry is 2.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	260	
1	B	260	

2 Entry composition ⓘ

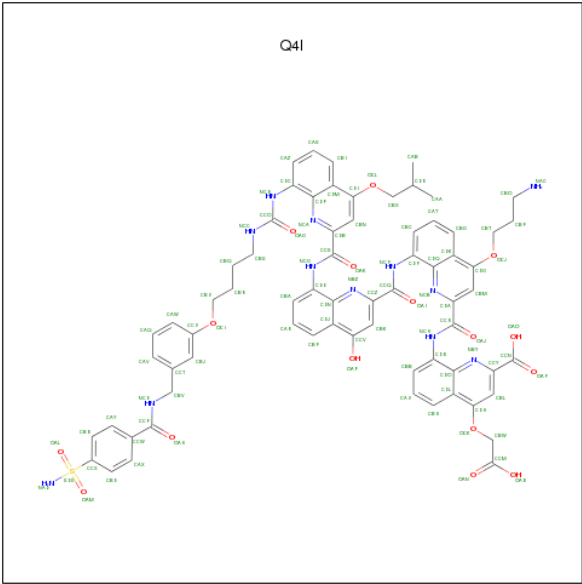
There are 4 unique types of molecules in this entry. The entry contains 4276 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Carbonic anhydrase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	S	0	0	0
			1981	1276	335	368	2			
1	B	257	Total	C	N	O	S	0	0	0
			1946	1253	324	367	2			

- Molecule 2 is 8-({[4-(3-AMINOPROPOXY)-8-({[4-HYDROXY-8-({[4-(2-METHYLPROPOXY)-8-({[4-(3-{{[(4-SULFAMOYLBENZOYL)AMINO]METHYL}PHENOXY)BUTYL]CARBAMOYL}AMINO)QUINOLIN-2-YL]CARBOONYL}AMINO)QUINOLIN-2-YL]CARBOONYL}AMINO)QUINOLIN-2-YL]CARBOONYL}AMINO)-4-(CARBOXYMETHOXY)QUINOLINE-2-CARBOXYLICACID (three-letter code: Q4I) (formula: C₆₈H₆₄N₁₂O₁₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			91	65	11	14	1		
2	B	1	Total	C	N	O	S	0	0
			97	68	12	16	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total 4	Zn 4	0	0
3	A	9	Total 9	Zn 9	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	84	Total 84	O 84	0	0
4	B	64	Total 64	O 64	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

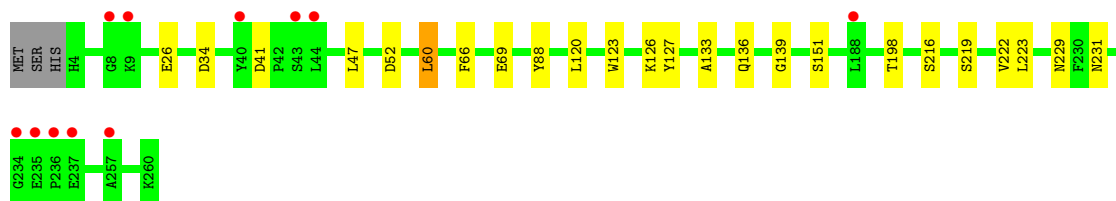
- Molecule 1: Carbonic anhydrase 2

Chain A: 



- Molecule 1: Carbonic anhydrase 2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	45.96Å 84.17Å 76.70Å 90.00° 97.18° 90.00°	Depositor
Resolution (Å)	41.50 – 2.15 41.47 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.7 (41.50-2.15) 99.8 (41.47-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.191 , 0.253 0.197 , 0.254	Depositor DCC
R_{free} test set	1582 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 31627 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4276	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, Q4I

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.83	1/2042 (0.0%)	0.84	1/2786 (0.0%)
1	B	0.66	0/2004	0.77	1/2742 (0.0%)
All	All	0.75	1/4046 (0.0%)	0.81	2/5528 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	7	TYR	CE1-CZ	5.99	1.46	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	34	ASP	CB-CG-OD1	9.12	126.51	118.30
1	B	34	ASP	CB-CG-OD1	6.71	124.33	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1981	0	1845	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1946	0	1774	11	0
2	A	91	0	54	1	0
2	B	97	0	62	3	0
3	A	9	0	0	0	0
3	B	4	0	0	0	0
4	A	84	0	0	1	0
4	B	64	0	0	2	0
All	All	4276	0	3735	24	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (24) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:69:GLU:OE2	4:B:457:HOH:O	2.18	0.57
1:A:127:TYR:CE1	1:A:136:GLN:HG3	2.44	0.52
1:A:88:TYR:HA	1:A:123:TRP:O	2.10	0.52
1:A:169:LYS:NZ	4:A:449:HOH:O	2.30	0.51
1:B:88:TYR:HA	1:B:123:TRP:O	2.12	0.50
1:B:133:ALA:O	1:B:139:GLY:HA3	2.12	0.50
1:B:219:SER:O	1:B:223:LEU:HD12	2.13	0.48
2:A:301:Q4I:OAG	2:A:301:Q4I:H32	2.15	0.47
1:B:219:SER:HA	1:B:222:VAL:HG12	1.97	0.47
1:A:133:ALA:O	1:A:139:GLY:HA3	2.16	0.46
1:A:41:ASP:C	1:A:41:ASP:OD1	2.54	0.46
1:B:41:ASP:C	1:B:41:ASP:OD1	2.54	0.46
1:B:127:TYR:CE1	1:B:136:GLN:HG3	2.52	0.45
1:A:229:ASN:HB3	1:A:231:ASN:OD1	2.17	0.44
2:B:301:Q4I:H32	2:B:301:Q4I:OAG	2.18	0.44
2:B:301:Q4I:H12	2:B:301:Q4I:OAI	2.19	0.43
1:B:60:LEU:O	1:B:66:PHE:HA	2.19	0.43
1:B:229:ASN:HB3	1:B:231:ASN:OD1	2.19	0.43
1:B:198:THR:OG1	2:B:301:Q4I:NAD	2.53	0.41
1:A:219:SER:HA	1:A:222:VAL:HG12	2.03	0.41
1:A:60:LEU:O	1:A:66:PHE:HA	2.21	0.41
1:B:26:GLU:HG2	4:B:456:HOH:O	2.21	0.40
1:A:127:TYR:CZ	1:A:136:GLN:HG3	2.57	0.40
1:A:34:ASP:HB3	1:A:37:THR:OG1	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	255/260 (98%)	243 (95%)	11 (4%)	1 (0%)	43	39
1	B	255/260 (98%)	243 (95%)	11 (4%)	1 (0%)	43	39
All	All	510/520 (98%)	486 (95%)	22 (4%)	2 (0%)	43	39

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	LYS
1	B	126	LYS

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	200/225 (89%)	192 (96%)	8 (4%)	42	41
1	B	192/225 (85%)	186 (97%)	6 (3%)	52	54
All	All	392/450 (87%)	378 (96%)	14 (4%)	47	47

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	LEU
1	A	52	ASP
1	A	57	LEU
1	A	60	LEU
1	A	120	LEU

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Mol	Chain	Res	Type
1	A	148	LYS
1	A	151	SER
1	A	216	SER
1	B	47	LEU
1	B	52	ASP
1	B	60	LEU
1	B	120	LEU
1	B	151	SER
1	B	216	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	53	GLN
1	A	229	ASN
1	B	4	HIS
1	B	229	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 13 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	Q4I	A	301	3	99,100,106	1.75	13 (13%)	142,143,150	1.54	22 (15%)
2	Q4I	B	301	3	106,106,106	2.07	20 (18%)	150,150,150	2.03	30 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	Q4I	A	301	3	-	0/64/64/71	0/10/10/10
2	Q4I	B	301	3	-	0/71/71/71	0/10/10/10

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	Q4I	CCX-SDS	-9.63	1.62	1.77
2	A	301	Q4I	CCY-CCN	-9.32	1.38	1.50
2	B	301	Q4I	CBV-NCD	7.61	1.61	1.46
2	B	301	Q4I	SDS-NAD	6.86	1.75	1.60
2	B	301	Q4I	CCY-CCN	-6.19	1.42	1.50
2	A	301	Q4I	CDA-CCR	-5.20	1.38	1.50
2	B	301	Q4I	CDB-CCS	-4.94	1.38	1.50
2	A	301	Q4I	CDB-CCS	-4.94	1.38	1.50
2	B	301	Q4I	CCW-CCP	-4.62	1.40	1.50
2	A	301	Q4I	OAM-SDS	3.63	1.50	1.43
2	A	301	Q4I	OAP-CCV	3.53	1.45	1.35
2	A	301	Q4I	OAL-SDS	3.51	1.50	1.43
2	B	301	Q4I	CDC-NCE	-3.48	1.31	1.41
2	B	301	Q4I	CCZ-CCQ	-3.44	1.42	1.50
2	A	301	Q4I	CAX-CCW	3.21	1.44	1.39
2	A	301	Q4I	CCW-CCP	-3.09	1.43	1.50
2	B	301	Q4I	CDA-CCR	-3.06	1.43	1.50
2	B	301	Q4I	CBF-CDJ	-2.87	1.36	1.42
2	B	301	Q4I	CDO-NBY	-2.76	1.31	1.37
2	B	301	Q4I	CDF-NCF	-2.71	1.33	1.41
2	A	301	Q4I	CCZ-CCQ	-2.69	1.44	1.50
2	B	301	Q4I	CDK-CDQ	-2.46	1.35	1.42
2	A	301	Q4I	CDI-CDM	-2.36	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	Q4I	CDM-CDP	-2.33	1.36	1.42
2	B	301	Q4I	OAL-SDS	2.31	1.47	1.43
2	B	301	Q4I	CBH-CDL	-2.28	1.37	1.42
2	B	301	Q4I	CCP-NCD	-2.24	1.29	1.33
2	A	301	Q4I	CDH-CDL	-2.21	1.36	1.42
2	A	301	Q4I	CBV-CCT	-2.17	1.46	1.51
2	B	301	Q4I	CCV-CDJ	-2.09	1.39	1.43
2	B	301	Q4I	CAX-CBD	2.04	1.42	1.38
2	B	301	Q4I	CDH-CDL	-2.02	1.36	1.42
2	A	301	Q4I	CBF-CDJ	-2.02	1.38	1.42

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	Q4I	OAM-SDS-NAD	-10.67	90.14	107.32
2	B	301	Q4I	CBV-NCD-CCP	-7.90	100.78	121.79
2	A	301	Q4I	OAL-SDS-CCX	-6.89	98.21	107.40
2	B	301	Q4I	OAO-CCN-CCY	6.11	128.47	114.60
2	B	301	Q4I	OAH-CCP-NCD	-5.13	111.64	122.55
2	B	301	Q4I	CCN-CCY-NBY	4.88	123.56	116.42
2	A	301	Q4I	CCN-CCY-NBY	4.72	123.31	116.42
2	B	301	Q4I	CCW-CCP-NCD	4.59	127.74	117.13
2	B	301	Q4I	CCU-CBJ-CCT	4.44	124.34	119.71
2	B	301	Q4I	CAX-CBD-CCX	4.26	124.34	119.47
2	A	301	Q4I	CBK-CCV-CDJ	-4.17	115.99	120.47
2	B	301	Q4I	CBD-CAX-CCW	-3.99	116.08	120.76
2	B	301	Q4I	OAP-CCV-CDJ	3.60	121.49	116.20
2	B	301	Q4I	CBM-CDA-NCB	-3.55	116.32	124.81
2	B	301	Q4I	OAL-SDS-NAD	3.55	113.03	107.32
2	B	301	Q4I	OAO-CCN-OAF	-3.49	115.50	123.36
2	A	301	Q4I	CCV-CDJ-CDN	3.45	122.05	118.19
2	B	301	Q4I	CDA-NCB-CDQ	3.37	125.19	117.03
2	A	301	Q4I	OAM-SDS-NAD	3.26	112.56	107.32
2	A	301	Q4I	CCV-CBK-CCZ	3.20	121.34	118.96
2	A	301	Q4I	OAO-CCN-CCY	3.12	121.67	114.60
2	A	301	Q4I	CBH-CDL-CDH	-3.04	114.92	122.51
2	B	301	Q4I	CBD-CCX-CBE	-3.03	115.66	120.40
2	A	301	Q4I	CBG-CDK-CDG	3.00	122.23	119.41
2	A	301	Q4I	CCX-SDS-NAD	2.98	113.07	108.51
2	B	301	Q4I	CBL-CCY-NBY	-2.90	117.88	124.81
2	B	301	Q4I	OAL-SDS-OAM	2.88	123.95	118.71
2	B	301	Q4I	CAY-CCW-CCP	-2.83	111.80	120.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	Q4I	CBL-CCY-CCN	-2.77	114.68	119.57
2	B	301	Q4I	CCV-CBK-CCZ	2.71	120.97	118.96
2	A	301	Q4I	CAU-CBH-CDL	-2.67	117.03	120.88
2	B	301	Q4I	OAF-CCN-CCY	-2.65	115.64	121.16
2	B	301	Q4I	OAM-SDS-CCX	2.65	110.94	107.40
2	B	301	Q4I	CBK-CCZ-NBZ	-2.63	118.53	124.81
2	B	301	Q4I	CAX-CCW-CCP	2.61	128.68	120.57
2	B	301	Q4I	CCY-NBY-CDO	2.52	123.15	117.03
2	A	301	Q4I	CBN-CDB-NCA	-2.43	119.00	124.81
2	B	301	Q4I	CBQ-CBS-NCC	-2.41	105.02	112.20
2	A	301	Q4I	CDA-NCB-CDQ	2.39	122.82	117.03
2	B	301	Q4I	CCZ-NBZ-CDN	2.38	122.81	117.03
2	A	301	Q4I	CBK-CCZ-NBZ	-2.36	119.16	124.81
2	A	301	Q4I	OCJ-CDG-CDK	-2.35	116.26	119.57
2	A	301	Q4I	CBF-CDJ-CCV	-2.24	118.86	122.13
2	A	301	Q4I	OAO-CCN-OAF	-2.23	118.34	123.36
2	A	301	Q4I	CCU-CBJ-CCT	2.17	121.97	119.71
2	A	301	Q4I	OAK-CCS-NCG	2.13	128.45	123.72
2	B	301	Q4I	CAY-CBE-CCX	2.12	121.90	119.47
2	A	301	Q4I	CBC-CDF-NCF	2.12	129.07	123.06
2	A	301	Q4I	CAS-CBI-CDM	-2.12	117.83	120.88
2	B	301	Q4I	CBW-OCK-CDH	-2.10	114.67	116.85
2	B	301	Q4I	CDF-NCF-CCQ	-2.04	122.59	128.88
2	B	301	Q4I	CBB-CDE-NCH	2.03	128.82	123.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	257/260 (98%)	-0.19	0 100 100	29, 42, 62, 80	0
1	B	257/260 (98%)	0.35	11 (4%) 34 37	33, 55, 77, 96	0
All	All	514/520 (98%)	0.08	11 (2%) 60 65	29, 48, 73, 96	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	PRO	5.6
1	B	234	GLY	4.6
1	B	235	GLU	2.8
1	B	9	LYS	2.7
1	B	237	GLU	2.7
1	B	44	LEU	2.7
1	B	40	TYR	2.6
1	B	8	GLY	2.6
1	B	43	SER	2.5
1	B	188	LEU	2.5
1	B	257	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	Q4I	A	301	91/97	0.14	1.18	24,41,55,59	0
3	ZN	A	310	1/1	0.12	0.81	28,28,28,28	0
3	ZN	B	305	1/1	0.15	0.75	36,36,36,36	0
2	Q4I	B	301	97/97	0.13	0.32	26,36,60,78	0
3	ZN	A	304	1/1	0.12	0.32	63,63,63,63	1
3	ZN	A	302	1/1	0.11	-0.15	34,34,34,34	0
3	ZN	B	302	1/1	0.11	-0.46	51,51,51,51	0
3	ZN	A	303	1/1	0.15	-1.06	53,53,53,53	1
3	ZN	B	303	1/1	0.12	-1.17	79,79,79,79	1
3	ZN	A	306	1/1	0.08	-1.71	55,55,55,55	0
3	ZN	B	304	1/1	0.05	-3.33	86,86,86,86	0
3	ZN	A	309	1/1	0.03	-3.85	64,64,64,64	0
3	ZN	A	308	1/1	0.04	-4.81	75,75,75,75	0
3	ZN	A	307	1/1	0.06	-5.34	64,64,64,64	0
3	ZN	A	305	1/1	0.03	-8.70	71,71,71,71	1

6.5 Other polymers

There are no such residues in this entry.